

Data Path : Z:\HPCHEM1\BNA F\DATA\BF050117\
 Data File : BF094753.D
 Acq On : 1 May 2017 16:06
 Operator : SJ/MA
 Sample : I2927-01
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FK-GF-03-003-A

Quant Time: May 02 02:49:29 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF041717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 17 14:59:23 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	5.76	152	110911	20.00	ng	0.00
21) Naphthalene-d8	7.25	136	448839	20.00	ng	0.00
38) Acenaphthene-d10	9.39	164	177444	20.00	ng	0.00
63) Phenanthrene-d10	11.21	188	295486	20.00	ng	0.00
75) Chrysene-d12	14.47	240	226540	20.00	ng	0.00
86) Perylene-d12	16.09	264	185231	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	4.32	112	742533	111.07	ng	0.02
7) Phenol-d6	5.40	99	892630	113.26	ng	0.00
23) Nitrobenzene-d5	6.41	82	541922	74.55	ng	0.00
41) 2,4,6-Tribromophenol	10.37	330	150869	108.70	ng	0.00
44) 2-Fluorobiphenyl	8.58	172	933905	77.44	ng	0.00
78) Terphenyl-d14	13.19	244	582180	68.69	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	9.08	163	132868	10.05	ng	98
70) Phenanthrene	11.24	178	109898	6.60	ng	96
74) Fluoranthene	12.70	202	155322	9.01	ng	99
77) Pyrene	12.98	202	140905	8.90	ng	99
80) Benzo(a)anthracene	14.45	228	63792	4.84	ng	98
82) Chrysene	14.49	228	59687	4.96	ng	96
85) Indeno(1,2,3-cd)pyrene	17.38	276	25716	2.25	ng	96
87) Benzo(b)fluoranthene	15.69	252	71323m	6.29	ng	
88) Benzo(k)fluoranthene	15.72	252	23694m	2.21	ng	
89) Benzo(a)pyrene	16.04	252	49305	4.68	ng	# 93
91) Benzo(g,h,i)perylene	17.76	276	25276	2.84	ng	96

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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